

IN THE CLAIMS:

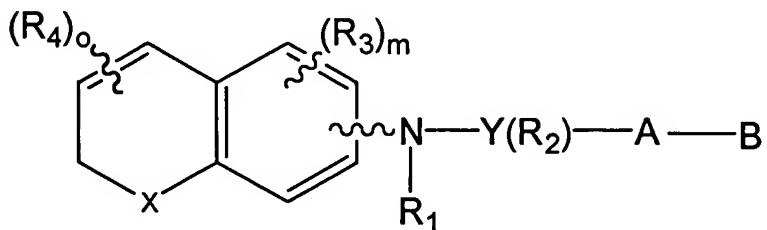
Cancel Claim 3.

Amend Claim 1 as set forth below in the Complete Listing of Pending

Claims:

COMPLETE LISTING OF PENDING CLAIMS

1. (Currently amended) A compound of the formula



where **X** is O, S, or C(R)₂;

R is H or alkyl of 1 to 6 carbons;

R₁ is H, alkyl of 1 to 10 carbons, alkenyl of 2 to 6 carbons; phenyl-C₁ - C₆ alkyl, or C₁ - C₆-alkylphenyl;

R₂ is H, alkyl of 1 to 6 carbons, F, Cl, Br, I, CF₃, fluoro substituted alkyl of 1 to 6 carbons, alkoxy of 1 to 6 carbons, or alkylthio of 1 to 6 carbons;

R₃ is independently alkyl of 1 to 6 carbons, F, Cl, Br, I, CF₃, fluoro substituted alkyl of 1 to 6 carbons, OH, SH, alkoxy of 1 to 10 carbons, fluoroalkoxy of 1 to 6 carbons, alkylthio of 1 to 6 carbons; benzyloxy C₁ - C₆ alkyl substituted benzyloxy, halogen substituted benzyloxy, phenoxy, C₁ - C₆ alkyl substituted phenoxy, or halogen substituted phenoxy;

R₄ is independently H, alkyl of 1 to 6 carbons, or F;

Y is a naphthyl group, or heteroaryl selected from a group consisting of pyridyl, thienyl, furyl, pyridazinyl, pyrimidinyl, pyrazinyl, thiazolyl, oxazolyl, imidazolyl and pyrazolyl group, said phenyl and heteroaryl groups pyridyl group being optionally substituted with one or two **R**₂ groups;

m is an integer having the values 0 to 3;

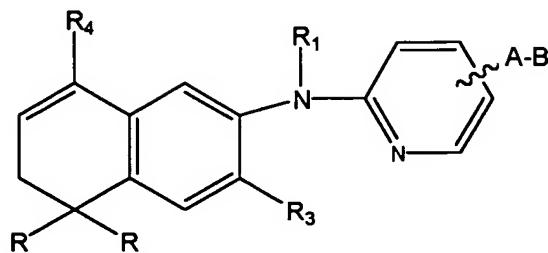
o is an integer having the values 0 to 4;

A is $(CH_2)_q$ where q is 0-5, lower branched chain alkyl having 3-6 carbons, cycloalkyl having 3-6 carbons, alkenyl having 2-6 carbons and 1 or 2 double bonds, alkynyl having 2-6 carbons and 1 or 2 triple bonds, and

B is hydrogen, COOH, COOR₈, CONR₉R₁₀, -CH₂OH, CH₂OR₁₁, CH₂OCOR₁₁, CHO, CH(OR₁₂)₂, CHOR₁₃O, -COR₇, CR₇(OR₁₂)₂, CR₇OR₁₃O, or tri-lower alkylsilyl, where R₇ is an alkyl, cycloalkyl or alkenyl group containing 1 to 5 carbons, R₈ is an alkyl group of 1 to 10 carbons or trimethylsilylalkyl where the alkyl group has 1 to 10 carbons, or a cycloalkyl group of 5 to 10 carbons, or R₈ is phenyl or lower alkylphenyl, R₉ and R₁₀ independently are hydrogen, an alkyl group of 1 to 10 carbons, or a cycloalkyl group of 5-10 carbons, or phenyl or lower alkylphenyl, R₁₁ is lower alkyl, phenyl or lower alkylphenyl, R₁₂ is lower alkyl, and R₁₃ is divalent alkyl radical of 2-5 carbons, or a pharmaceutically acceptable salt of said compound.

2. (original) A compound in accordance with Claim 1 where **X** is C(**R**)₂.
3. (cancelled)
4. (original) A compound in accordance with Claim 1 where **X** is S.
5. (original) A compound in accordance with Claim 1 where **X** is O.
6. (original) A compound in accordance with Claim 1 where the **A-B** group represents $(CH_2)_qCOOR_8$ or $(CH_2)_qCOOH$ where q is 0, or a pharmaceutically acceptable salt thereof.
7. (original) A compound in accordance with Claim 1 where **R**₁ is alkyl of 1 to 10 carbons or alkenyl of 2 to 6 carbons.
8. (original) A compound in accordance with Claim 1 where **R**₄ is independently H or alkyl of 1 to 6 carbons.
9. (original) A compound in accordance with Claim 1 where **R**₁ is alkyl of 1 to 10 carbons or alkenyl of 2 to 6 carbons, **R**₄ is independently H or alkyl of 1 to 6 carbons and the **A-B** group represents $(CH_2)_qCOOR_8$ or $(CH_2)_qCOOH$ where q is 0, or a pharmaceutically acceptable salt thereof.

10. (previously submitted) A compound that has the structure of formula (iii)



(iii)

where **R** is independently H or alkyl of 1 to 6 carbons;

R₁ is H or alkyl of 1 to 10 carbons or alkenyl of 2 to 6 carbons;

R₃ is independently alkyl of 1 to 6 carbons, F, Cl, Br, I, CF₃, fluoro substituted alkyl of 1 to 6 carbons, OH, SH, alkoxy of 1 to 10 carbons, fluoroalkoxy of 1 to 6 carbons, alkylthio of 1 to 6 carbons; benzyloxy, C₁ - C₆ alkyl substituted benzyloxy, halogen substituted benzyloxy, phenoxy, C₁ - C₆ alkyl substituted phenoxy, or halogen substituted phenoxy;

R₄ is H or alkyl of 1 to 6 carbons;

A is (CH₂)_q where q is 0-5, lower branched chain alkyl having 3-6 carbons, cycloalkyl having 3-6 carbons, alkenyl having 2-6 carbons and 1 or 2 double bonds, alkynyl having 2-6 carbons and 1 or 2 triple bonds, and

B is hydrogen, COOH, COOR₈, CONR₉R₁₀, -CH₂OH, CH₂OR₁₁, CH₂OCOR₁₁, CHO, CH(OR₁₂)₂, CHOR₁₃O, -COR₇, CR₇(OR₁₂)₂, CR₇OR₁₃O, or tri-lower alkylsilyl, where R₇ is an alkyl, cycloalkyl or alkenyl group containing 1 to 5 carbons, R₈ is an alkyl group of 1 to 10 carbons or trimethylsilylalkyl where the alkyl group has 1 to 10 carbons, or a cycloalkyl group of 5 to 10 carbons, or R₈ is phenyl or lower alkylphenyl, R₉ and R₁₀ independently are hydrogen, an alkyl group of 1 to 10 carbons, or a cycloalkyl group of 5-10 carbons, or phenyl or lower alkylphenyl, R₁₁ is lower alkyl, phenyl or lower alkylphenyl, R₁₂ is lower alkyl, and R₁₃ is divalent alkyl radical of 2-5 carbons, or a pharmaceutically

acceptable salt of said compound.

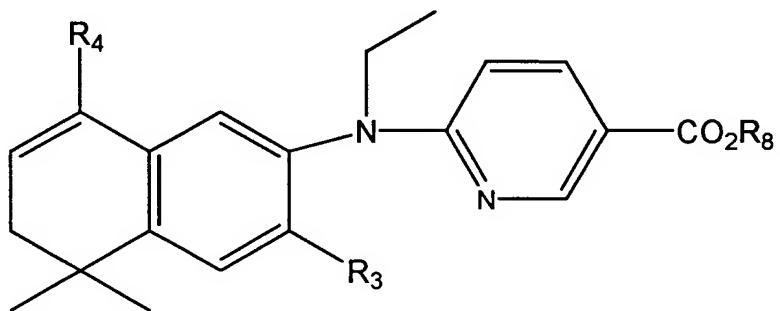
11 – 13 inclusive (canceled)

14. (original) A compound in accordance with Claim 10 where **R₄** and **R₁** both are alkyl.

15. (original) A compound in accordance with Claim 10 where the **A-B** group represents $(CH_2)_qCOOR_8$ or $(CH_2)_qCOOH$ where **q** is 0, or a pharmaceutically acceptable salt thereof.

16 – 27 inclusive (canceled)

28. (original) A compound of the formula



where **R₃** is H, or alkyl of 1 to 6 carbons;

R₄ is alkyl of 1 to 6 carbons, and

R₈ is H or alkyl of 1 to 6 carbons, or a pharmaceutically acceptable salt of said compound.

29. (original) A compound in accordance with Claim 28 where **R₄** is methyl, ethyl, *i*-propyl or *t*-butyl.

30. (original) A compound in accordance with Claim 28 where **R₃** is H, or *n*-butyloxy.

31. (original) A compound in accordance with Claim 28 where **R₈** is H or ethyl.

32. (original) A compound in accordance with Claim 28 where **R₄** is methyl, ethyl, *i*-propyl or *t*-butyl;

R₃ is H, or *n*-butyloxy, and **R₈** is H or ethyl.

33. (original) A compound in accordance with Claim 32 where **R₈** is H or ethyl.